	Type	L #	Hits	Search Text	DBs	Time Stamp
1	BRS	L3	75	((PQQ$ or pyrrolo$) near3 dehydrogenase)	USPAT; EPO; JPO; DERWENT	2002/01/13 15:29
2	BRS	L4	8	3 and (phthal$ or mal$ or succin$)	USPAT; EPO; JPO; DERWENT	2002/01/13 15:32

	Type	L #	Hits	Search Text	DBs	Time Stamp
1	BRS	L6	21	(sens$ or biosens$ or detect$ or measur$) and electrode$ and (PQQ$ or pyrrlo$) and glucon$	USPAT; EPO; JPO; DERWENT	2002/01/13 14:12
2	BRS	L7	2	(glucose near1 dehydrogenase) and (glucon$ near5 (stabiliz$ or additive))	USPAT; EPO; JPO; DERWENT	2002/01/13 14:15
3	BRS	L8	815	(glucose dehydrogenase).ti.	USPAT	2002/01/13 14:20
4	BRS	L9	9	(glucose near1 dehydrogenase).ti.	USPAT	2002/01/13 14:24
5	BRS	L10	19	(glucose near1 dehydrogenase).ti. or (glucose near1 dehydrogenase).ab.	USPAT	2002/01/13 14:24
6	BRS	L11	4	10 and (gluconic or gluconate)	USPAT; EPO; JPO; DERWENT	2002/01/13 14:25

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=> d his full

     (FILE 'HOME' ENTERED AT 13:50:39 ON 13 JAN 2002)

     FILE 'CAPLUS' ENTERED AT 13:51:04 ON 13 JAN 2002
L1           212 SEA ((GLUCOSE DEHYDROGENASE) AND (GLUCON?))/AB
L2            10 SEA L1 AND (STABIL?)/AB
                 D 1-10 ALL
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=> d his full

     (FILE 'HOME' ENTERED AT 12:58:26 ON 13 JAN 2002)

     FILE 'REGISTRY' ENTERED AT 12:58:42 ON 13 JAN 2002
L1             2 SEA GLUCONIC ACID/CN
                 D ALL

     FILE 'CAPLUS' ENTERED AT 12:59:44 ON 13 JAN 2002
L2           126 SEA (SENSOR OR BIOSENSOR OR DETECT?) AND ELECTRODE? AND
                 (GLUCON?)
L3            35 SEA L2 AND (PQQ? OR PYRROLO? OR DEHYDROGENASE)
                 D 1-35 ALL

     FILE 'STNGUIDE' ENTERED AT 13:02:36 ON 13 JAN 2002
L4             0 SEA (SENS? OR MEASUR?) AND ELECTRODE? AND (GLUCON?)

     FILE 'CAPLUS' ENTERED AT 13:12:45 ON 13 JAN 2002
L5           191 SEA (SENS? OR MEASUR?) AND ELECTRODE? AND (GLUCON?)
L6            32 SEA L5 AND (PQQ? OR PYRROLO? OR DEHYDROGENASE)
                 D 10 ALL
L7             4 SEA L6 NOT L3
                 D 1-4 ALL

     FILE 'STNGUIDE' ENTERED AT 13:13:32 ON 13 JAN 2002
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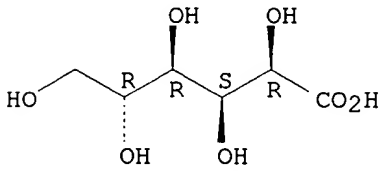
L3 ANSWER 1 OF 2 REGISTRY COPYRIGHT 2002 ACS
RN 526-95-4 REGISTRY
CN D-Gluconic acid (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN Gluconic acid, D- (8CI)
OTHER NAMES:
CN **Gluconic acid**
AR 133-42-6
FS STEREOSEARCH
MF C6 H12 O7
CI COM
LC STN Files: AGRICOLA, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS, BIOTECHNO, CA, CABA, CANCERLIT, CAPLUS, CASREACT, CEN, CHEMCATS, CHEMINFORMRX, CHEMLIST, CIN, CSCHEM, DIOGENES, EMBASE, GMELIN*, HODOC*, HSDB*, IFICDB, IFIPAT, IFIUDB, IPA, MEDLINE, MRCK*, MSDS-OHS, NIOSHTIC, PDLCOM*, PIRA, PROMT, RTECS*, TOXCENTER, TULSA, ULIDAT, USAN, USPAT2, USPATFULL
(*File contains numerically searchable property data)
Other Sources: DSL**, EINECS**, TSCA**
(**Enter CHEMLIST File for up-to-date regulatory information)

Absolute stereochemistry.



Calculated Properties (CALC)

CODE	PROPERTY	VALUE	CONDITION	NOTE
HD	H donors	6		ACD (1)
HAC	H acceptors	7		ACD (1)
MW	Molecular Weight	196.16		ACD (1)
LOGP	logP	-3.175+/-0.852		ACD (1)
LOGD	logD	-3.18	pH 1	ACD (1)
LOGD	logD	-3.91	pH 4	ACD (1)
LOGD	logD	-6.69	pH 7	ACD (1)
LOGD	logD	-7.17	pH 8	ACD (1)
LOGD	logD	-7.27	pH 10	ACD (1)
PKA	pKa	3.35+/-0.35	Most Acidic	ACD (1)
SLB.MOL	Molar Solubility	>=1 mol/L	pH 1	ACD (1)
SLB.MOL	Molar Solubility	>=1 mol/L	pH 4	ACD (1)
SLB.MOL	Molar Solubility	>=1 mol/L	pH 7	ACD (1)
SLB.MOL	Molar Solubility	>=1 mol/L	pH 8	ACD (1)
SLB.MOL	Molar Solubility	>=1 mol/L	pH 10	ACD (1)

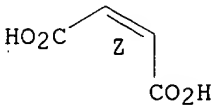
(1) Calculated using Advanced Chemistry Development (ACD) Software Solaris V4.67 ((C) 1994-2002 ACD)

4012 REFERENCES IN FILE CA (1967 TO DATE)
546 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
4023 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1

L4 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2002 ACS
RN 110-16-7 REGISTRY
CN 2-Butenedioic acid (2Z)- (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN 2-Butenedioic acid (Z)-
CN **Maleic acid (8CI)**
OTHER NAMES:
CN 2-Butenedioic acid, (Z)-
CN cis-1,2-Ethylenedicarboxylic acid
CN cis-2-Butenedioic acid
CN cis-Butenedioic acid
CN Maleinic acid
CN Malezid CM
CN Scotchbond Multipurpose Etchant
CN Toxilic acid
FS STEREOSEARCH
MF C4 H4 O4
CI COM
LC STN Files: AGRICOLA, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS,
BIOTECHNO, CA, CABA, CANCERLIT, CAOLD, CAPLUS, CASREACT, CBNB, CEN,
CHEMCATS, CHEMINFORMRX, CHEMLIST, CIN, CSCHEM, CSNB, DDFU, DETHERM*,
DIPPR*, DRUGU, EMBASE, ENCOMPLIT, ENCOMPLIT2, ENCOMPPAT, ENCOMPPAT2,
GMELIN*, HODOC*, HSDB*, IFICDB, IFIPAT, IFIUDB, IPA, MEDLINE, MRCK*,
MSDS-OHS, NAPRALERT, NIOSHTIC, PDLCOM*, PIRA, PROMT, RTECS*, SPECINFO,
TOXCENTER, TULSA, USPAT2, USPATFULL
(*File contains numerically searchable property data)
Other Sources: DSL**, EINECS**, TSCA**
(**Enter CHEMLIST File for up-to-date regulatory information)

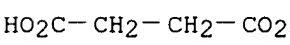
Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

9345 REFERENCES IN FILE CA (1967 TO DATE)
2696 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
9365 REFERENCES IN FILE CAPLUS (1967 TO DATE)
5 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

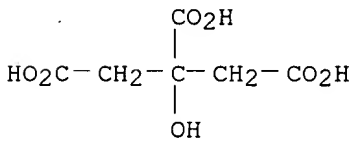
L5 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2002 ACS
RN 110-15-6 REGISTRY
CN Butanedioic acid (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN **Succinic acid (8CI)**
OTHER NAMES:
CN 1,2-Ethanedicarboxylic acid
CN 1,4-Butanedioic acid
CN A 12084
CN Amber acid
CN Asuccin
CN Dihydrofumaric acid
CN Katasuccin
CN Wormwood acid
FS 3D CONCORD
MF C4 H6 O4
CI COM
LC STN Files: AGRICOLA, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS,
BIOTECHNO, CA, CABA, CANCERLIT, CAOLD, CAPLUS, CASREACT, CBNB, CEN,
CHEMCATS, CHEMINFORMRX, CHEMLIST, CIN, CSCHEM, DDFU, DETHERM*, DIOGENES,
DIPPR*, DRUGU, EMBASE, ENCOMPLIT, ENCOMPLIT2, ENCOMPPAT, ENCOMPPAT2,
GMELIN*, HODOC*, HSDB*, IFICDB, IFIPAT, IFIUDB, IPA, MEDLINE, MRCK*,
MSDS-OHS, NAPRALERT, NIOSHTIC, PDLCOM*, PIRA, PROMT, RTECS*, SPECINFO,
SYNTHLINE, TOXCENTER, TULSA, ULIDAT, USPAT2, USPATFULL, VETU, VTB
(*File contains numerically searchable property data)
Other Sources: DSL**, EINECS**, TSCA**
(**Enter CHEMLIST File for up-to-date regulatory information)



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

17334 REFERENCES IN FILE CA (1967 TO DATE)
1954 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
17363 REFERENCES IN FILE CAPLUS (1967 TO DATE)
9 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

L6 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2002 ACS
RN 77-92-9 REGISTRY
CN 1,2,3-Propanetricarboxylic acid, 2-hydroxy- (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN **Citric acid (8CI)**
OTHER NAMES:
CN 2-Hydroxy-1,2,3-propanetricarboxylic acid
CN 3-Carboxy-3-hydroxypentane-1,5-dioic acid
CN Aciletten
CN Chemfill
CN Citretten
CN Citro
CN F 0001 (polycarboxylic acid)
CN Hydrocerol A
CN Uro-trainer
FS 3D CONCORD
DR 12262-73-6, 43136-35-2, 245654-34-6
MF C6 H8 O7
CI COM
LC STN Files: ADISNEWS, AGRICOLA, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS, BIOTECHNO, CA, CABA, CANCERLIT, CAOLD, CAPLUS, CASREACT, CBNB, CEN, CHEMCATS, CHEMINFORMRX, CHEMLIST, CIN, CSCHEM, CSNB, DDFU, DETHERM*, DIOGENES, DIPPR*, DRUGU, EMBASE, ENCOMPLIT, ENCOMPLIT2, ENCOMPPAT, ENCOMPPAT2, GMELIN*, HODOC*, HSDB*, IFICDB, IFIPAT, IFIUDB, IPA, MEDLINE, MRCK*, MSDS-OHS, NAPRALERT, NIOSHTIC, PDLCOM*, PHARMASEARCH, PIRA, PROMT, RTECS*, SPECINFO, SYNTHLINE, TOXCENTER, TULSA, USAN, USPAT2, USPATFULL, VETU, VTB
(*File contains numerically searchable property data)
Other Sources: DSL**, EINECS**, TSCA**
(**Enter CHEMLIST File for up-to-date regulatory information)



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

36859 REFERENCES IN FILE CA (1967 TO DATE)
2437 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
36945 REFERENCES IN FILE CAPLUS (1967 TO DATE)
9 REFERENCES IN FILE CAOLD (PRIOR TO 1967)